

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112318\
 Data File : BF110948.D
 Acq On : 23 Nov 2018 17:07
 Operator : JU/SJ
 Sample : J5999-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSSI-1115-S-AB-2

Manual Integrations
 APPROVED

MOHAMMAD
 11/26/2018 3:12:07 PM

Quant Time: Nov 23 21:24:27 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	68992	20.00	ng	-0.01
21) Naphthalene-d8	8.22	136	275145	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	111934	20.00	ng	-0.02
64) Phenanthrene-d10	11.47	188	189744	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	167490	20.00	ng	-0.01
87) Perylene-d12	15.66	264	129668	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	400674	94.33	ng	0.00
7) Phenol-d6	6.56	99	492374	90.65	ng	-0.01
23) Nitrobenzene-d5	7.50	82	311325	65.16	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	85991	76.24	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	355612	45.37	ng	-0.02
79) Terphenyl-d14	13.05	244	279160	31.21	ng	-0.02
Target Compounds						
10) Phenol	6.57	94	26621	4.227	ng	Qvalue # 77
50) Dimethylphthalate	9.69	163	36422	4.361	ng	99
71) Phenanthrene	11.50	178	77537	7.627	ng	99
72) Anthracene	11.56	178	24293	2.369	ng	97
75) Fluoranthene	12.69	202	152540	14.967	ng	99
78) Pyrene	12.93	202	137975	10.699	ng	98
81) Benzo(a)anthracene	14.12	228	83936	8.384	ng	98
83) Chrysene	14.15	228	74757	7.838	ng	98
86) Indeno(1,2,3-cd)pyrene	17.25	276	29656	4.333	ng	95
88) Benzo(b)fluoranthene	15.20	252	86278m	11.122	ng	
89) Benzo(k)fluoranthene	15.23	252	32649m	4.259	ng	
90) Benzo(a)pyrene	15.59	252	58065	8.239	ng	97
92) Benzo(a,h,i)perylene	17.73	276	28791	4.865	ng	# 95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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